

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021517\
 Data File : BF092952.D
 Acq On : 15 Feb 2017 15:51
 Operator : SJ/MA
 Sample : PB96964BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB96964BS

Manual Integrations
 APPROVED

mohammad
 2/16/2017 11:40:07 AM

Quant Time: Feb 15 16:22:12 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 13 18:04:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	140715	20.00	ng	-0.02
21) Naphthalene-d8	8.16	136	624205	20.00	ng	-0.01
38) Acenaphthene-d10	9.92	164	333732	20.00	ng	-0.01
63) Phenanthrene-d10	11.39	188	523628	20.00	ng	-0.01
75) Chrysene-d12	14.03	240	478309	20.00	ng	-0.02
86) Perylene-d12	15.46	264	386889	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.48	112	1179979	143.87	ng	0.00
7) Phenol-d6	6.52	99	1588128	150.49	ng	-0.01
23) Nitrobenzene-d5	7.44	82	918308	81.93	ng	-0.02
41) 2,4,6-Tribromophenol	10.70	330	314576	130.33	ng	-0.01
44) 2-Fluorobiphenyl	9.23	172	1615132	87.68	ng	-0.01
78) Terphenyl-d14	12.98	244	1400601	68.80	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.52	88	161644	36.47	ng	86
3) Pyridine	3.25	79	457152	40.57	ng	86
4) n-Nitrosodimethylamine	3.21	42	225870	42.69	ng	85
6) Aniline	6.53	93	231560	16.09	ng	# 70
8) 2-Chlorophenol	6.66	128	436271	48.21	ng	98
9) Benzaldehyde	6.42	77	99140	13.11	ng	92
10) Phenol	6.53	94	528395	42.80	ng	# 76
11) bis(2-Chloroethyl)ether	6.61	93	445714	45.23	ng	98
12) 1,3-Dichlorobenzene	6.81	146	457622	43.18	ng	# 92
13) 1,4-Dichlorobenzene	6.89	146	485450	45.26	ng	95
14) 1,2-Dichlorobenzene	7.04	146	412728	40.24	ng	96
15) Benzyl Alcohol	7.01	79	351774	45.02	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.15	45	719280	42.01	ng	93
17) 2-Methylphenol	7.14	107	384848	49.58	ng	95
18) Hexachloroethane	7.38	117	154686	42.24	ng	91
19) n-Nitroso-di-n-propylamine	7.29	70	293728	43.72	ng	96
20) 3+4-Methylphenols	7.29	107	432671	46.62	ng	90
22) Acetophenone	7.29	105	566987	39.78	ng	# 92
24) Nitrobenzene	7.46	77	502560	43.66	ng	93
25) Isophorone	7.70	82	892038	42.71	ng	99
26) 2-Nitrophenol	7.78	139	248136	45.35	ng	# 77
27) 2,4-Dimethylphenol	7.81	122	389414	40.53	ng	97
28) bis(2-Chloroethoxy)methane	7.90	93	533190	38.54	ng	99
29) 2,4-Dichlorophenol	8.02	162	402896	44.96	ng	93
30) 1,2,4-Trichlorobenzene	8.10	180	393073	40.70	ng	97
31) Naphthalene	8.18	128	1264841	39.97	ng	99
32) Benzoic acid	7.95	122	279231	50.80	ng	98
33) 4-Chloroaniline	8.22	127	89278	7.19	ng	95
34) Hexachlorobutadiene	8.29	225	204513	38.49	ng	99
35) Caprolactam	8.60	113	119842	42.51	ng	# 28
36) 4-Chloro-3-methylphenol	8.72	107	412932	44.20	ng	97
37) 2-Methylnaphthalene	8.86	142	808456	39.79	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.04	216	374492	39.98	ng	99
40) Hexachlorocyclopentadiene	9.02	237	342420	81.01	ng	95

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42) 2,4,6-Trichlorophenol	9.15	196	276425	44.91	ng	91
43) 2,4,5-Trichlorophenol	9.20	196	275682	40.87	ng #	89
45) 1,1'-Biphenyl	9.33	154	980579	40.58	ng	97
46) 2-Chloronaphthalene	9.36	162	769147	40.69	ng	96
47) 2-Nitroaniline	9.46	65	286603	45.09	ng	92
48) Acenaphthylene	9.78	152	1325533	40.59	ng	98
49) Dimethylphthalate	9.64	163	1019360	45.35	ng	100
50) 2,6-Dinitrotoluene	9.70	165	193991	39.96	ng #	77
51) Acenaphthene	9.95	154	866103	44.36	ng	96
52) 3-Nitroaniline	9.87	138	83424	15.02	ng #	79
53) 2,4-Dinitrophenol	9.98	184	195324	79.15	ng	94
54) Dibenzofuran	10.12	168	1152317	44.12	ng	90
55) 4-Nitrophenol	10.04	139	298833	74.68	ng	95
56) 2,4-Dinitrotoluene	10.11	165	304737	47.15	ng #	77
57) Fluorene	10.46	166	847659	42.19	ng	97
58) 2,3,4,6-Tetrachlorophenol	10.24	232	215110	47.81	ng	99
59) Diethylphthalate	10.34	149	944647	44.59	ng	96
60) 4-Chlorophenyl-phenylether	10.45	204	409289	42.40	ng #	81
61) 4-Nitroaniline	10.49	138	220267	38.71	ng	92
62) Azobenzene	10.61	77	1074176	49.08	ng	94
64) 4,6-Dinitro-2-methylphenol	10.51	198	129809	41.13	ng #	39
65) n-Nitrosodiphenylamine	10.57	169	737357	44.08	ng	99
66) 4-Bromophenyl-phenylether	10.94	248	247591	45.26	ng #	80
67) Hexachlorobenzene	11.00	284	226515	41.90	ng #	76
68) Atrazine	11.10	200	260040	48.44	ng	94
69) Pentachlorophenol	11.21	266	277150	97.57	ng	98
70) Phenanthrene	11.42	178	1295213	43.17	ng	98
71) Anthracene	11.47	178	1205560	40.02	ng	99
72) Carbazole	11.63	167	1227589	42.63	ng	98
73) Di-n-butylphthalate	11.95	149	1465482	47.05	ng	98
74) Fluoranthene	12.60	202	1208745	39.06	ng	98
76) Benzidine	12.73	184	268979	13.20	ng	96
77) Pyrene	12.83	202	1202533	33.60	ng	100
79) Butylbenzylphthalate	13.45	149	579383	39.86	ng	95
80) Benzo(a)anthracene	14.02	228	1074435	36.73	ng	99
81) 3,3'-Dichlorobenzidine	13.99	252	228532	21.92	ng #	95
82) Chrysene	14.05	228	955090	34.87	ng	99
83) Bis(2-ethylhexyl)phthalate	14.01	149	822317	41.37	ng #	97
84) Di-n-octyl phthalate	14.61	149	1420544	43.88	ng	99
85) Indeno(1,2,3-cd)pyrene	16.88	276	881580	41.55	ng #	94
87) Benzo(b)fluoranthene	15.05	252	1223887m	48.06	ng	
88) Benzo(k)fluoranthene	15.08	252	737632m	33.55	ng	
89) Benzo(a)pyrene	15.40	252	930265	42.23	ng	97
90) Dibenzo(a,h)anthracene	16.89	278	730507	41.03	ng #	94
91) Benzo(g,h,i)perylene	17.31	276	746612	41.51	ng #	94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

